



Microarticle

Tellurium coincidence Doppler broadening spectroscopy study

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ABSTRACT

The Coincidence Doppler broadening (CDB) reference curve for pure Tellurium (Te) has been obtained stabilising a point of departure for future Positron Annihilation Spectroscopy (PAS) studies of alloys that include Te among its components to determine the influence of this element in the open volume defects and its environment.

Introduction

Te is a p-type semiconductor widely used because its numerous properties such as photoconductivity [1], nonlinear optical responses [2], and high-performance thermoelectric behavior [3] among others. These properties allow Te to be involved in a wide different uses such as solar cells [4]; biological theranostic (therapeutic or diagnostic) applications [5]; or in the new developing high-entropy alloys to improve its mechanical properties [6]. A complete characterization of Te and Te-based materials is required for any further development. PAS allows studying the vacancy-type defects that strongly influence the mechanical properties of metallic alloys [7] and semiconductors [8], nevertheless reference values and curves for pure Te are required and they are not present in literature [9,10]. In this study CDB reference curve corresponding to pure Te has been obtained.

Experimental procedure

The midway position between the two HP Ge detectors (placed face to face separated by 40 cm to each other, and configured in 1 μ s timing coincidence) was used as target position for the samples that sandwiched a 0.5 MBq ^{22}Na source sealed in kapton. The 24 spectra taken showed no evidence of electronic shift and had a total number of events around 3×10^7 in the 512×512 coincidence matrix. The obtained cumulative spectra after selecting a strip in the diagonal of the matrix had 1×10^7 counts. The selection window was defined as $2m_0c^2 - 1.6 \text{ keV} < E_1 + E_2 < 2m_0c^2 + 1.6 \text{ keV}$, where E_1 and E_2 stand for the energies of the pair of annihilating photons, m_0 the electron rest mass and c the light speed. Using a bin width of $2.5 \times 10^{-3} m_0$ improved the shape of the curve for the resulting data, and then normalized for comparison purposes. As a reference for the above data, 99.999% purity annealed (3 h 400 °C) Al samples (10 mm \times 10 mm, 2 mm thick) analogous CDB spectrum was used in order to highlight the signature points of the sample's spectra.

Results

Normalized to pure Al CDB curve for Te (99.999% purity, annealed at 250 °C during 2 h, 7 mm \times 7 mm, 5 mm thick) shown in Fig. 1 a) shows a first major peak at $p_L = 15.2(3) \times 10^{-3} m_0c$ and a second peak at $p_L = 21.2(3) \times 10^{-3} m_0c$ followed by a constant decrease in the high-momentum region. Fe and Cu obtained CDB curves displayed in Fig. 1 a) match theoretical calculations [11] confirming the experimental results of the setup. The momentum position of the maximum of the first peak can be placed in the distribution of [10] and compared to the other elements of the $36 < Z < 54$ region that partially share the electronic momentum distribution ($[\text{Kr}]5s^1$ to $[\text{Kr}]5s^24d^{10}5p^6$). The obtained point, located in the end of the Z-grouped elements, is located on top of the other data following the same tendency of the previous Z region ($18 < Z < 36$) where all the elements showed a quick incensement in the distribution.

Conclusions

CDB spectroscopy has been used to characterize high-purity Te. The obtained Te CDB curve has been validated by the reference plot “position of the first peak of the curve” versus “atomic number Z” [10] obtaining the corresponding new point at $p_L = 15.2(3) \times 10^{-3} m_0c$. This curve will be used as reference in further studies involving Te-based materials.

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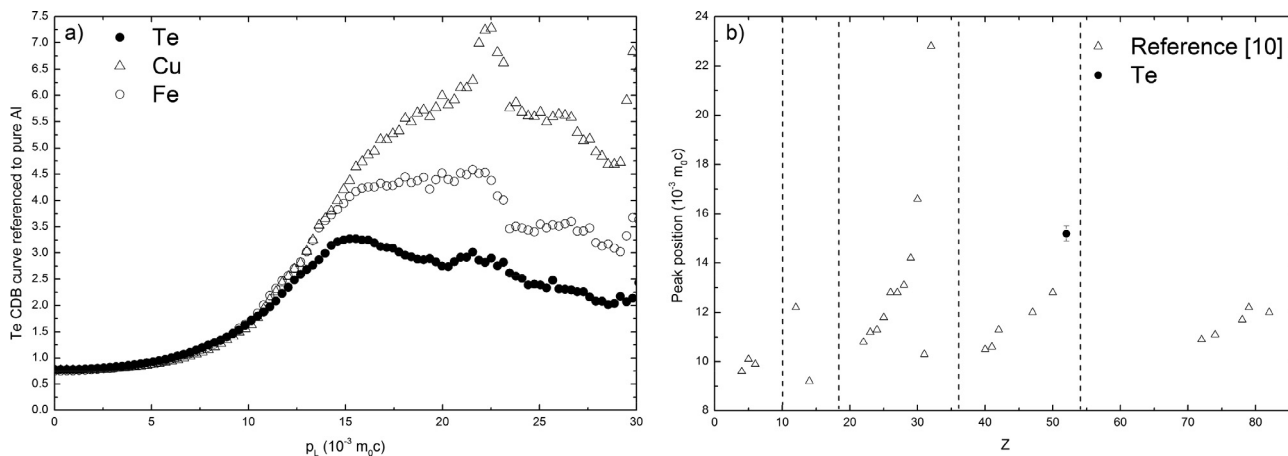


Fig. 1. a) Te CDB curve referenced to pure Al. Fe and Cu CDB curves added for comparison purposes. b) CDB first peak position as a function of atomic number Z . Solid dot correspond to reference [10] and empty dot correspond to this work.

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